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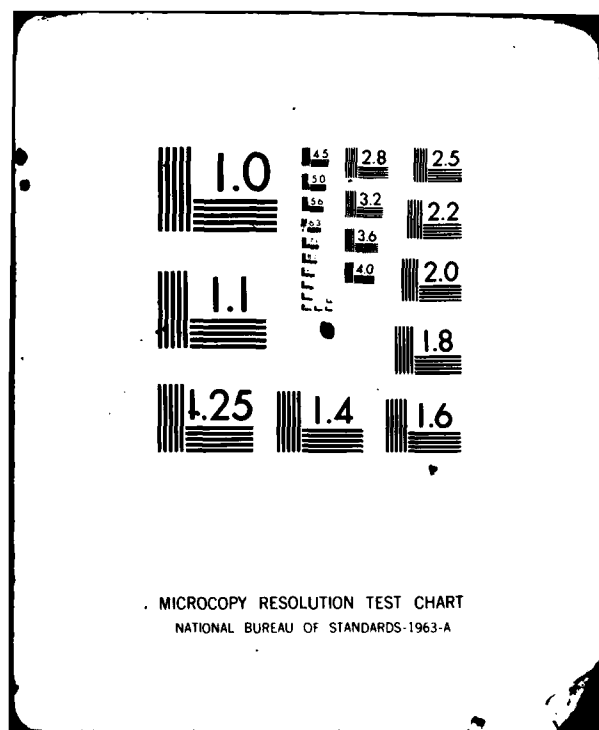
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THE CHOLESKY DECOMPOSITION ALGORITHM AND ARMA MODELING

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The Cholesky Decomposition Algorithm and ARMA Modeling

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Summary

The history of the use of the Cholesky decomposition in the estimation and prediction of autoregressive moving average time series is reviewed and a modification of Ansley's (1979) estimation method is suggested. This modification is motivated by the results of Newton and Pagano (1981) in prediction theory and results in a significant reduction in the calculations required to evaluate the exact likelihood. It is also shown how the modified method can be used to obtain predictors and prediction variances.

Some key words: Autoregressive moving average model; prediction; Likelihood estimation; Cholesky decomposition.



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1. Introduction

Let $\{\epsilon(t), t = 0, \pm 1, \dots\}$ be a white noise time series of zero mean uncorrelated random variables having variance σ^2 . Then the time series $\{Y(t), t = 0, \pm 1, \dots\}$ satisfying

$$\sum_{j=0}^p \alpha(j)Y(t-j) = \sum_{k=0}^q \beta(k)\epsilon(t-k), \quad \alpha(0) = \beta(0) = 1,$$

is called an autoregressive moving average process of order p, q (ARMA (p,q)). If $q = 0$, $Y(\cdot)$ is called an autoregressive process of order p , AR(p), while if $p = 0$ then $Y(\cdot)$ is a moving average process of order q , MA(q). Modern time series analysis has found the ARMA process extremely useful in modeling and predicting future values of a time series. However, the computational tasks involved in obtaining efficient estimators of the parameters $\alpha(\cdot)$, $\beta(\cdot)$, and σ^2 and in obtaining predictors of future values of $Y(\cdot)$ are extremely difficult and have led to the development of a wide variety of methods for accomplishing these tasks. These methods approach the problem from many points of view including 1) modified Newton Raphson likelihood maximization (Hannan (1969), Parzen (1971), Akaike (1973), for example), 2) nonlinear least squares (Box and Jenkins (1970), for example), 3) Kalman filtering techniques (Akaike (1974), Harvey and Phillips (1979), Jones (1980), Pearlman (1980), for example), 4) the use of the Cholesky matrix triangularization technique (Whittle (1963), Pagano and Parzen (1973), Pagano (1976), Pagano (1973), Phadke and Kedem (1978), Ansley (1979), Newton (1980), and Newton and Pagano (1981), for example), and others.

The purpose of this paper is to review the history of the Cholesky algorithm as applied to ARMA models and to show how one can easily modify the estimation procedure of Ansley (1979), which is currently regarded as the most computationally efficient exact likelihood estimation technique for at least the case where $p \geq q$ or $q < 5$ (see Pearlman (1980)), to obtain a significant reduction in the number of computations required.

Section 2 describes the ARMA process estimation and prediction problems and reviews the use of the Cholesky algorithm, while the proposed modification is given in section 3.

2. The Estimation and Prediction Problems

Given a realization $\underline{Y}_T = (Y(1), \dots, Y(T))^T$ from an ARMA(p,q) process, one often desires estimators of the $\alpha(\cdot)$, $\beta(\cdot)$, and σ^2 as well as memory-t, horizon-h, minimum mean square error linear predictors $Y(t+h|t)$ and prediction variances $\sigma_{t,h}^2 = E\{Y(t+h) - Y(t+h|t)\}^2$ of $Y(t+h)$ given $\underline{Y}_t = (Y(1), \dots, Y(t))^T$ for a variety of values h_1, \dots, h_2 and t_1, \dots, t_2 of t and h .

The usual procedure is to obtain a Gaussian likelihood identification $\hat{\alpha}(\cdot)$, $\hat{\beta}(\cdot)$, and $\hat{\sigma}^2$ as the values of the parameters maximizing the Gaussian likelihood function

$$L(\underline{\alpha}, \underline{\beta}, \sigma^2 | \underline{Y}_T) = (2\pi)^{-T/2} |\Gamma_{Y,T}|^{-1} \exp(-\underline{Y}_T^T \Gamma_{Y,T}^{-1} \underline{Y}_T)$$

where $\Gamma_{Y,t} = \text{Cov}(\underline{Y}_t) = (E(Y(j)Y(k))) = R_Y(|j-k|)$, $j, k = 1, \dots, t$. The Gaussian assumption is often used also so that $Y(t+h|t)$ and $\sigma_{t,h}^2$, which in general are given by $Y(t+h|t) = \lambda_{t,h}^T \underline{Y}_t$ and $\sigma_{t,h}^2 = R_Y(0) - \underline{r}_{t,h}^T \Gamma_{Y,t}^{-1} \underline{r}_{t,h}$ where $\Gamma_{t,h} \lambda_{t,h} = \underline{r}_{t,h} = (R_Y(t+h-1), \dots, R_Y(h))^T$, can be obtained as the conditional mean and variance of $Y(t+h)$ given \underline{Y}_t .

With the advent of efficient numerical algorithms for maximizing nonlinear functions that require a user to only supply a routine that evaluates the function for specified values of its arguments (see Dennis and More (1977), for example), the emphasis on finding the identification of $\alpha(\cdot)$, $\beta(\cdot)$, and σ^2 has been on finding efficient algorithms for evaluating L , or equivalently a more easily evaluated function having maxima at the same values of its parameters. The major efforts in this regard have been to apply the Kalman filter algorithm and the Cholesky decomposition algorithm. Each of these methods were originally used in ARMA modeling as a means of obtaining predictors and prediction variances.

2.1 The Cholesky Decomposition

Let A_n be a symmetric $(n \times n)$ matrix. Then A_n is positive definite if and only if it can be decomposed as the product $A_n = M_{A,n} M_{A,n}^T$ (Cholesky decomposition) or $A_n = L_{A,n} D_{A,n} L_{A,n}^T$ (modified Cholesky decomposition) where $M_{A,n}$ is an $(n \times n)$ lower triangular matrix with positive diagonal elements and $L_{A,n}$ is a unit lower triangular matrix while $D_{A,n}$ is a diagonal matrix containing the squares of the diagonal elements of $M_{A,n}$. Further these decompositions are unique and nested, i.e., for example, $L_{A,K}$ is the lower triangular matrix in the decomposition of the $K \times K$ principal minor of A_n , $K = 1, \dots, n$. Thus the $(j,k)^{th}$ element of $L_{A,K}$ can be denoted $L_{A,j,k}$ for $j, k \leq K$.

2.2 The Cholesky Decomposition and Prediction

In his seminal work on time series prediction, Whittle (1963) suggested that such a triangularization would be useful in finding $Y(t+h|t)$ and $\sigma_{t,h}^2$. Pagano and Parzen (1973) and Pagano (1976) showed that for an $MA(q)$ process

$$Y(t+h|t) = \begin{cases} \sum_{k=h}^q L_{Y,t+h,t+h-k} e(t+h-k), & h=0, \dots, q \\ 0 & , h>q \end{cases}$$

$$\sigma_{t,h}^2 = \sum_{k=0}^{h-1} L_{Y,t+h,t+h-k}^2 D_{Y,t+h-k,t+h-k}$$

$$L_{Y,K,K-j} = 0 \quad , j>q$$

$$\lim_{K \rightarrow \infty} L_{Y,K,K-j} = \beta(j) \quad , j = 1, \dots, q,$$

$$\lim_{K \rightarrow \infty} D_{Y,K,K} = \sigma^2,$$

where $e(1) = Y(1)$ and $e(j) = Y(j) - \sum_{k=1}^{\min(j-1,q)} L_{Y,j,j-k} e(j-k)$, $j>2$ is the j th element of $L_{Y,K}^{-1} Y_K$ for any $K \geq j$. Thus one need only calculate the successive rows of L_Y and the corresponding $e(\cdot)$ until convergence of its $q+1$ nonzero elements to the $\beta(\cdot)$ is achieved, at which point one replaces $L_{Y,K,K-j}$ and $D_{Y,K,K}$ by $\beta(j)$ and σ^2 respectively in the above formulas. Newton (1980) discusses the extension of this algorithm to the multivariate MA(q) process.

Pagano and Parzen (1973) further suggested that this procedure could be extended to the ARMA (p,q) case by applying the autoregressive filter $\alpha(\cdot)$ to $Y(p+1), \dots, Y(T)$ and using the MA(q) algorithm combined with an AR(p) prediction algorithm.

Finally Newton and Pagano (1981) show that for a general purely nondeterministic covariance stationary time series $Y(\cdot)$

$$Y(t+h|t) = \sum_{k=h}^{t+h-1} L_{Y,t+h,t+h-k} e(t+h-k)$$

$$\sigma_{t,h}^2 = \sum_{k=0}^{h-1} L_{Y,t+h,t+h-k}^2 D_{Y,t+h-k,t+h-k}$$

$$\lim_{K \rightarrow \infty} L_{Y,K,K-j} = \beta_{\infty}(j) \quad , \quad j \geq 0$$

$$\lim_{K \rightarrow \infty} D_{Y,K,K} = \sigma_{\infty}^2 \quad ,$$

where the $\beta_{\infty}(\cdot)$ and σ_{∞}^2 are the coefficients and the noise variance of the $MA(\infty)$ representation of $Y(\cdot)$ and $e(1) = Y(1)$, while

$$e(j) = Y(j) - \sum_{k=1}^{j-1} L_{Y,j,k} e(k) \quad , \quad j \geq 2 \quad .$$

Applying this general theory to the ARMA (p,q) process and further factoring L_Y into $L_Y = L_Z L_X$ where $Z(\cdot)$ is an AR(p) process with coefficients $\alpha(\cdot)$ and noise variance σ^2 and $\Gamma_{X,K} = L_{Z,K}^{-1} \Gamma_{Y,K} L_{Z,K}^{-T} = L_{X,K} D_{X,K} L_{X,K}^T$ has last $K-q$ rows and columns being the covariance matrix of an MA(q) process, Newton and Pagano obtain

$$Y(t+h|t) = X(t+h|t) - \sum_{j=1}^p \alpha(j) Y(t+h-j|t) \quad ,$$

$$\sigma_{t,h}^2 = \sum_{k=0}^{h-1} D_{X,t+h-k,t+h-k} \left\{ \sum_{\ell=t+h-k}^{t+h} L_{Z,t+h,\ell} L_{X,\ell,t+h-k} \right\}^2 \quad ,$$

$$Y(t+h-j|t) = Y(t+h-j) \quad \text{if } j \geq h \quad ,$$

$$X(t+h|t) = \begin{cases} \sum_{k=h}^q L_{X,t+h,t+h-k} e(t+h-k), & h = 1, \dots, q \\ 0, & h > q \end{cases} \quad ,$$

$e(j)$ is the j^{th} element of $L_{X,K}^{-1} X_K$, $K \geq j$, and $X_K = L_{Z,K}^{-1} Y_K$. Further,

$$\lim_{K \rightarrow \infty} L_{X,K,K-j} = \beta(j), \quad j = 1, \dots, q,$$

$$\lim_{K \rightarrow \infty} D_{X,K,K} = \sigma^2.$$

A number of other results are given by Newton and Pagano which greatly reduce the amount of storage and computations necessary for obtaining L_X and L_Z . We note that this method is different than that given by Pagano and Parzen (1973) in that instead of leaving the initial values $Y(1), \dots, Y(p)$ untransformed, the Newton and Pagano algorithm forms the time series $X(\cdot)$ by $X_K = L_{Z,K}^{-1} Y_K$, i.e.

$$X(j) = Y(j), \quad j = 1$$

$$\sum_{k=0}^{j-1} \alpha_{j-1}(k) Y(j-k), \quad j = 2, \dots, p$$

$$\sum_{k=0}^p \alpha(k) Y(j-k), \quad j > p$$

where the $\alpha_j(k)$ are easily obtained by performing the Levinson (1974) recursion for $j = p-1, \dots, 1$, with $\alpha_{p+1}(k) = \alpha(k)$:

$$\alpha_j(i) = \frac{\alpha_{j+1}(i) - \alpha_{j+1}(j+1)\alpha_{j+1}(j+1-i)}{1 - \alpha_{j+1}^2(j+1)}, \quad i = 1, \dots, j < p$$

2.3 The Cholesky Decomposition and Parameter Estimation

Pagano (1973) discusses the use of the Cholesky decomposition in solving the Yule Walker equations for an AR(p) process. Phadke and Kedem (1978) and Ansley (1979) apply the Cholesky decomposition to the problem

of evaluating the likelihood function of the MA(q) and ARMA(p,q) processes respectively, essentially using the same transformation and decomposition that Pagano and Parzen (1973) used for obtaining predictors. We briefly describe Ansley's method and show in the next section how it can be improved by incorporating the theory of section 2.2.

Define a new time series, with $m = \max(p, q+1)$,

$$W(t) = Y(t), \quad t = 1, \dots, m$$

$$\sum_{j=0}^p \alpha(j)Y(t-j), \quad t = m+1, \dots, T.$$

Thus in matrix notation this is $\underline{W}_T = \underline{M}_{Y,T} \underline{Y}_T$ where $\underline{M}_{Y,T}$ is a (TXT) unit lower triangular matrix and $\underline{\Gamma}_{W,T} = \text{Cov}(\underline{W}_T) = \underline{M}_{Y,T} \underline{\Gamma}_{Y,T} \underline{M}_{Y,T}^T$, which is a banded matrix of bandwidth q, i.e., there are q nonzero subdiagonals, from row $m+1$ on. Thus finding the Cholesky decomposition $\underline{\Gamma}_{W,T} = \underline{M}_{W,T} \underline{M}_{W,T}^T$ and the new random vector $\underline{e}_T = \underline{M}_{W,T}^{-1} \underline{W}_T$ can be done recursively only having to compute q elements in each row of $\underline{M}_{W,T}$. Then $\underline{e}_T = \underline{M}_{W,T}^{-1} \underline{M}_{Y,T} \underline{Y}_T$ and thus

$$\underline{e}_T \sim N_T(0, \underline{M}_{W,T}^{-1} \underline{M}_{Y,T} \underline{\Gamma}_{Y,T} \underline{M}_{Y,T}^T \underline{M}_{W,T}^{-T} = \underline{I}_T).$$

Thus maximizing L is equivalent to maximizing

$$\rho(\underline{\alpha}, \underline{\beta}, \sigma^2 | \underline{e}_T) = (2\pi)^{-T/2} |\underline{M}_{W,T}|^{-1} \exp\left(-\sum_{t=1}^T \frac{e^2(t)}{2\sigma^2}\right).$$

3. Using the Prediction Results in Ansley's Method

Thus it is clear that Ansley's method consists of factoring the lower triangular matrix in the Cholesky decomposition of $\underline{\Gamma}_{Y,T}$ into the product of the two easily calculated lower triangular matrices $\underline{M}_{Y,T}$ and $\underline{M}_{W,T}$. But from section 2.2 it is easily seen that if this factorization is done as

$L_{Y,T} = L_{Z,T} L_{X,T}$ one can call upon the results of Newton and Pagano to halt calculation of successive rows of $L_{X,T}$ once their elements have converged to $\beta(1), \dots, \beta(q)$. The experience of the present authors is that the speed of this convergence, which is a function of how close the zeros of $\sum_{k=0}^q \beta(k)z^k$ are to the unit circle, is such that a significant reduction in computations is obtained, particularly if T is large relative to q . Further one can then use the prediction theory of section 2.2 to obtain desired predictors and prediction variances.

We note that section 2.2 can also be used to prove that the elements of the rows of Ansley's $M_{W,T}$ also converge to the $\beta(\cdot)$ and our experience is that the rate of convergence is virtually identical with the above method.

Thus either factorization could be used but it seems more natural to factor $L_{Y,T}$ into an exact AR(p) covariance matrix factor $L_{Z,T}$ and an exact MA(q) (except for the first $\max(p,q+1)$ rows and columns) covariance matrix factor $L_{X,T}$ and use the methods of section 2.2 to find the likelihood, the predictors, and the prediction variances.

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